

=> fil reg

FILE 'REGISTRY' ENTERED AT 16:22:31 ON 27 NOV 2007

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STRUCTURE FILE UPDATES: 26 NOV 2007 HIGHEST RN 955995-34-3

DICTIONARY FILE UPDATES: 26 NOV 2007 HIGHEST RN 955995-34-3

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TSCA INFORMATION NOW CURRENT THROUGH June 29, 2007

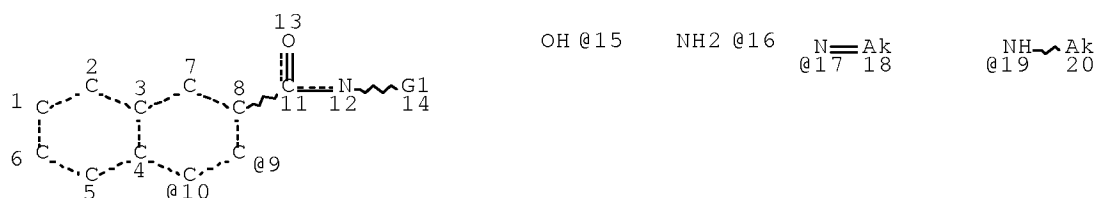
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<http://www.cas.org/support/stngen/stdoc/properties.html>

=> d que stat l19

L17 STR



VAR G1=16/17/19

VPA 15-9/10 U

NODE ATTRIBUTES:

CONNECT IS E2 RC AT 12

CONNECT IS E1 RC AT 18

CONNECT IS E1 RC AT 20

DEFAULT MLEVEL IS ATOM

GGCAT IS SAT AT 18

GGCAT IS SAT AT 20

DEFAULT ECLEVEL IS LIMITED

GRAPH ATTRIBUTES:

RSPEC I

NUMBER OF NODES IS 20

STEREO ATTRIBUTES: NONE

L19 45 SEA FILE=REGISTRY SSS FUL L17

100.0% PROCESSED 37921 ITERATIONS

45 ANSWERS

SEARCH TIME: 00.00.01

=> d his nofile

(FILE 'HOME' ENTERED AT 15:43:48 ON 27 NOV 2007)

FILE 'HCAPLUS' ENTERED AT 15:43:57 ON 27 NOV 2007

L1 1 SEA ABB=ON PLU=ON US2007157846/PN
SEL RN

FILE 'REGISTRY' ENTERED AT 15:44:29 ON 27 NOV 2007

L2 13 SEA ABB=ON PLU=ON (214417-91-1/BI OR 25014-41-9/BI OR
25067-34-9/BI OR 5341-58-2/BI OR 7439-89-6/BI OR
7440-50-8/BI OR 7440-66-6/BI OR 79-10-7/BI OR 79-41-4/BI
OR 9002-86-2/BI OR 9002-88-4/BI OR 9003-20-7/BI OR
9003-53-6/BI)

D SCA

L3 1 SEA ABB=ON PLU=ON L2 AND C11 H10 N2 O2/MF
D SCA

L4 1 SEA ABB=ON PLU=ON L2 AND C17 H20 N2 O2/MF
D SCA

FILE 'HCAPLUS' ENTERED AT 15:50:00 ON 27 NOV 2007

L5 110 SEA ABB=ON PLU=ON L3

L6 27 SEA ABB=ON PLU=ON L4

FILE 'STNGUIDE' ENTERED AT 15:51:14 ON 27 NOV 2007

L7 QUE ABB=ON PLU=ON ANTIRUST? OR RUSTPROOF? OR (INHIBIT?
OR PREVENT? OR PROHIBIT? OR ANTI) (A) (WEAR? OR CORRO? OR
OXID? OR RUST?)

L8 1 SEA ABB=ON PLU=ON (L5 OR L6) AND L7

L9 QUE ABB=ON PLU=ON MIX? OR BLEND? OR ADMIX? OR COMMIX?
OR IMMIX? OR INTERMIX? OR COMPOSIT? OR FORMULAT? OR
COMBINAT?

L10 54 SEA ABB=ON PLU=ON (L5 OR L6) AND L9

FILE 'LREGISTRY' ENTERED AT 16:03:44 ON 27 NOV 2007

L11 STR

L12 STR

FILE 'REGISTRY' ENTERED AT 16:07:32 ON 27 NOV 2007

L13 0 SEA SSS SAM L11

L14 STR 214417-91-1

L15 STR L11

L16 50 SEA SSS SAM L15

FILE 'LREGISTRY' ENTERED AT 16:10:14 ON 27 NOV 2007

L17 STR L15

FILE 'REGISTRY' ENTERED AT 16:14:58 ON 27 NOV 2007

L18 3 SEA SSS SAM L17

L19 45 SEA SSS FUL L17

L20 2 SEA ABB=ON PLU=ON L2 AND L19
SAV L19 GRE261/A

FILE 'HCAPLUS' ENTERED AT 16:15:50 ON 27 NOV 2007

L21 153 SEA ABB=ON PLU=ON L19

L22 2 SEA ABB=ON PLU=ON L21 AND L7

```
=> fil hcap
FILE 'HCAPLUS' ENTERED AT 16:22:43 ON 27 NOV 2007
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FILE COVERS 1907 - 27 Nov 2007 VOL 147 ISS 23
FILE LAST UPDATED: 26 Nov 2007 (20071126/ED)
```

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This file contains CAS Registry Numbers for easy and accurate substance identification.

```
=> d l3 ide
YOU HAVE REQUESTED DATA FROM FILE 'REGISTRY' - CONTINUE? (Y)/N:n
```

```
=> fil reg
FILE 'REGISTRY' ENTERED AT 16:22:52 ON 27 NOV 2007
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```
STRUCTURE FILE UPDATES: 26 NOV 2007 HIGHEST RN 955995-34-3
DICTIONARY FILE UPDATES: 26 NOV 2007 HIGHEST RN 955995-34-3
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TSCA INFORMATION NOW CURRENT THROUGH June 29, 2007

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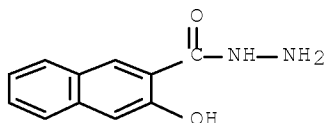
REGISTRY includes numerically searchable data for experimental and predicted properties as well as tags indicating availability of experimental property data in the original document. For information on property searching in REGISTRY, refer to:

<http://www.cas.org/support/stngen/stdoc/properties.html>

```
=> d ide l3
```

```
L3 ANSWER 1 OF 1 REGISTRY COPYRIGHT 2007 ACS on STN
RN 5341-53-2 REGISTRY
```

ED Entered STN: 16 Nov 1984
 CN 2-Naphthalenecarboxylic acid, 3-hydroxy-, hydrazide (CA INDEX NAME)
 OTHER CA INDEX NAMES:
 CN 2-Naphthoic acid, 3-hydroxy-, hydrazide (6CI, 7CI, 8CI)
 OTHER NAMES:
 CN 2-Hydroxy-3-naphthoic acid hydrazide
 CN 2-Hydroxy-3-naphthoylhydrazine
 CN 3-Hydroxy-2-naphthoic acid hydrazide
 CN 3-Hydroxy-2-naphthoylhydrazine
 CN 3-Hydroxy-naphthalene-2-carboxylic acid hydrazide
 CN NSC 2117
 CN NSC 49198
 MF C11 H10 N2 O2
 CI COM
 LC STN Files: BEILSTEIN*, BIOSIS, CA, CAOLD, CAPLUS, CASREACT,
 CHEMCATS, CHEMLIST, CSCHEM, IFICDB, IFIPAT, IFIUDB, MEDLINE,
 SPECINFO, TOXCENTER, USPAT2, USPATFULL, USPATOLD
 (*File contains numerically searchable property data)
 Other Sources: EINECS**, NDSL**, TSCA**
 (**Enter CHEMLIST File for up-to-date regulatory information)

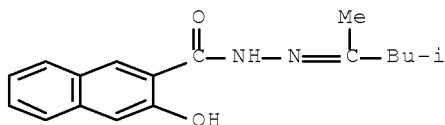


PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

110 REFERENCES IN FILE CA (1907 TO DATE)
 1 REFERENCES TO NON-SPECIFIC DERIVATIVES IN FILE CA
 110 REFERENCES IN FILE CAPLUS (1907 TO DATE)
 9 REFERENCES IN FILE CAOLD (PRIOR TO 1967)

=> d ide l4

L4 ANSWER 1 OF 1 REGISTRY COPYRIGHT 2007 ACS on STN
 RN 214417-91-1 REGISTRY
 ED Entered STN: 18 Nov 1998
 CN 2-Naphthalenecarboxylic acid, 3-hydroxy-, (1,3-dimethylbutylidene)hydrazide (9CI) (CA INDEX NAME)
 OTHER NAMES:
 CN BMH
 CN BMH (hydrazide)
 MF C17 H20 N2 O2
 SR CA
 LC STN Files: CA, CAPLUS, CHEMCATS, CHEMLIST, USPAT2, USPATFULL



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

27 REFERENCES IN FILE CA (1907 TO DATE)
 1 REFERENCES TO NON-SPECIFIC DERIVATIVES IN FILE CA
 27 REFERENCES IN FILE CAPLUS (1907 TO DATE)

=> fil hcap
 FILE 'HCAPLUS' ENTERED AT 16:23:06 ON 27 NOV 2007
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FILE COVERS 1907 - 27 Nov 2007 VOL 147 ISS 23
 FILE LAST UPDATED: 26 Nov 2007 (20071126/ED)

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=> d l22 ibib abs hitstr hitind 1-2

L22 ANSWER 1 OF 2 HCAPLUS COPYRIGHT 2007 ACS on STN
 ACCESSION NUMBER: 2005:903042 HCAPLUS Full-text
 DOCUMENT NUMBER: 143:233682
 TITLE: Corrosion inhibitor
 INVENTOR(S): Sonogi, Ken; Nabeshima, Akihiro
 PATENT ASSIGNEE(S): Otsuka Chemical Co., Ltd., Japan
 SOURCE: PCT Int. Appl., 19 pp.
 CODEN: PIXXD2
 DOCUMENT TYPE: Patent
 LANGUAGE: Japanese
 FAMILY ACC. NUM. COUNT: 1
 PATENT INFORMATION:

PATENT NO.	KIND	DATE	APPLICATION NO.	DATE
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WO 2005078157	A1	20050825	WO 2005-JP2823	

200502
 16

W: AE, AG, AL, AM, AT, AU, AZ, BA, BB, BG, BR, BW, BY, BZ, CA,
 CH, CN, CO, CR, CU, CZ, DE, DK, DM, DZ, EC, EE, EG, ES, FI,
 GB, GD, GE, GH, GM, HR, HU, ID, IL, IN, IS, JP, KE, KG, KP,
 KR, KZ, LC, LK, LR, LS, LT, LU, LV, MA, MD, MG, MK, MN, MW,
 MX, MZ, NA, NI, NO, NZ, OM, PG, PH, PL, PT, RO, RU, SC, SD,

SE, SG, SK, SL, SY, TJ, TM, TN, TR, TT, TZ, UA, UG, US, UZ,
 VC, VN, YU, ZA, ZM, ZW
 RW: BW, GH, GM, KE, LS, MW, MZ, NA, SD, SL, SZ, TZ, UG, ZM, ZW,
 AM, AZ, BY, KG, KZ, MD, RU, TJ, TM, AT, BE, BG, CH, CY, CZ,
 DE, DK, EE, ES, FI, FR, GB, GR, HU, IE, IS, IT, LT, LU, MC,
 NL, PL, PT, RO, SE, SI, SK, TR, BF, BJ, CF, CG, CI, CM, GA,
 GN, GQ, GW, ML, MR, NE, SN, TD, TG

TW 278444 B 20070411 TW 2005-94104280

200502
 15

EP 1717350 A1 20061102 EP 2005-710526

200502
 16

R: AT, BE, CH, DE, DK, ES, FR, GB, GR, IT, LI, LU, NL, SE, MC,
 PT, IE, SI, LT, FI, RO, CY, TR, BG, CZ, EE, HU, PL, SK, IS

CN 1922341 A 20070228 CN 2005-80005120

200502
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IN 2006DN04600 A 20070824 IN 2006-DN4600

200608
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US 2007157846 A1 20070712 US 2006-589261

200608
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PRIORITY APPLN. INFO.:

JP 2004-37782

A

200402
 16

WO 2005-JP2823

W

200502
 16

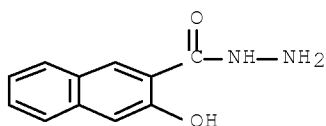
AB A corrosion inhibitor contains
 3-hydroxy-2-naphthoic acid hydrazide and/or 3-hydroxy-2-naphthoic acid (1,3-
 dimethylbutylidene)hydrazide and salts thereof as an active constituent. Also
 disclosed are a corrosion inhibition method and a corrosion
 inhibition resin composition using such a corrosion inhibitor. The corrosion
 inhibitor, method, and resin composition are suitable for protecting Fe, Cu,
 and Zn from corrosion.

IT 5341-58-2, 3-Hydroxy-2-naphthoic acid hydrazide
 214417-91-1

RL: NUU (Other use, unclassified); USES (Uses)
 (corrosion inhibitor)

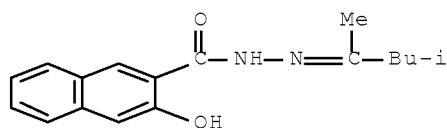
RN 5341-58-2 HCAPLUS

CN 2-Naphthalenecarboxylic acid, 3-hydroxy-, hydrazide (CA INDEX NAME)



RN 214417-91-1 HCAPLUS

CN 2-Naphthalenecarboxylic acid, 3-hydroxy-, (1,3-
 dimethylbutylidene)hydrazide (9CI) (CA INDEX NAME)



IC ICM C23F011-00
ICS B32B015-08; C07C243-38; C07C251-76; C09D005-08; C09D201-00

CC 56-10 (Nonferrous Metals and Alloys)

ST corrosion inhibitor copper iron zinc

IT Epoxy resins, uses
Polyamides, uses
Polycarbonates, uses
Polyesters, uses
Polyolefins
Polyurethanes, uses
RL: NUU (Other use, unclassified); USES (Uses)
(corrosion inhibitor composition)

IT Polyketones
RL: NUU (Other use, unclassified); USES (Uses)
(polyether-; corrosion inhibitor composition)

IT Polyethers, uses
RL: NUU (Other use, unclassified); USES (Uses)
(polyketone-; corrosion inhibitor composition)

IT 5341-58-2, 3-Hydroxy-2-naphthoic acid hydrazide
214417-91-1
RL: NUU (Other use, unclassified); USES (Uses)
(corrosion inhibitor)

IT 7439-89-6, Iron, uses 7440-50-8, Copper, uses 7440-66-6, Zinc, uses
RL: TEM (Technical or engineered material use); USES (Uses)
(corrosion inhibitor)

IT 79-10-7D, Acrylic acid, esters, polymers 79-41-4D, Methacrylic acid, esters, polymers 9002-86-2, S1001 9002-88-4, Sumikathene L-430 9003-20-7, Polyvinyl acetate 9003-53-6, Polystyrene 25014-41-9, Polyacrylonitrile 25067-34-9, EP-F101
RL: NUU (Other use, unclassified); USES (Uses)
(corrosion inhibitor composition)

REFERENCE COUNT: 5 THERE ARE 5 CITED REFERENCES AVAILABLE FOR THIS RECORD. ALL CITATIONS AVAILABLE IN THE RE FORMAT

L22 ANSWER 2 OF 2 HCAPLUS COPYRIGHT 2007 ACS on STN

ACCESSION NUMBER: 1962:38008 HCAPLUS Full-text

DOCUMENT NUMBER: 56:38008

ORIGINAL REFERENCE NO.: 56:7100g-i, 7101a-e

TITLE: Inhibitory effect of various hydrazides on monoamine oxidase in vitro and in vivo

AUTHOR(S): Szmuszkovicz, Jacob; Greig, Margaret E.

CORPORATE SOURCE: Upjohn Co., Kalamazoo, MI

SOURCE: Journal of Medicinal & Pharmaceutical Chemistry (1961), 4, 259-96
CODEN: JMPCAS; ISSN: 0095-9065

DOCUMENT TYPE: Journal

LANGUAGE: Unavailable

AB A series of aliphatic and aromatic hydrazides was prepared and tested as monoamine oxidase inhibitors. In vitro tests were done using the Bhagvat method in the Warburg apparatus The in vivo studies were done on several compds. by using oral activity in rats. Hydrazine hydrate (15 g.) in 25 ml.

EtOH was added to 75 g. diethyl oxalate in 25 ml. EtOH while holding the temperature at -15 to -25°, the suspension stirred at this temperature 10 min., allowed to warm to room temperature, stirred 1.25 hrs., filtered, the filtrate evaporated in vacuo at 30-5°, 75 ml. H₂O added, the solution extracted twice with Et₂O, the aqueous solution evaporated at 35° to remove Et₂O, 44 ml. Me₂CO added, allowed to stand 1 hr., evaporated in vacuo overnight at 25-30°, the solution freeze-dried, 250 ml. Me₂CO added, and the solution refluxed 2 hrs., and evaporated at 30° to give 37.7 g. of the isopropylidene derivative (I), m. 45-54°. I (37 g.) in 200 ml. EtOH was hydrogenated under 52 lb. H in the presence of 1 g. PtO₂. After 2/3 of the theoretical amount of H was absorbed 1 g. of catalyst was added and hydrogenation completed. The solution was evaporated to dryness at 40°, the residue dissolved in 50 ml. EtOH, added over 10 min. to 12.7 g. hydrazine hydrate in 100 ml. EtOH, the suspension stirred 1 hr., filtered, the solid washed with EtOH, refluxed with 2250 ml. EtOH, filtered, and allowed to crystallize to give 17.5 g. oxalic acid hydrazide 2-isopropylhydrazide, m. 178-81.5°. Oxalic acid hydrazide phenethylhydrazide, m. 153°, oxalic acid hydrazide 2-(α -benzyl)ethylhydrazide, m. 113-14°, oxalic acid bis(2-ethylhydrazide), m. 202.5-4.0°, oxalic acid bis(2-propylhydrazide), m. 193.5-4.5°, oxalic acid bis(2-isopropylhydrazide), m. 191.5-2.5°, oxalic acid 2-ethylhydrazide 2-isopropylhydrazide, m. 173-4°, oxalic acid bis(2-n-butylhydrazide), m. 157-8°, oxalic acid bis(2-isobutylhydrazide), m. 165-6°, oxalic acid bis[2-(1-methylpropyl)hydrazide], m. 135-6.5°, oxalic acid bis[2-(1-ethylpropyl)hydrazide], m. 118.5-19.5°, oxalic acid bis(2-benzylhydrazide), m. 165-7°, oxalic acid bis(phenethylhydrazide), m. 163-4°, oxalic acid bis(α -methylphenethylhydrazide), m. 159-60°, oxalic acid 2-isopropylhydrazide 2-(α -methylphenethyl)hydrazide, m. 150-3°, oxalic acid bis[2-(1-cyclopropylethyl)hydrazide], isomer I, m. 191-2°, isomer II, m. 163-4°, oxalic acid bis(2-cyclopentylhydrazide), m. 212-14°, oxalic acid 2-isopropylhydrazide 2-(α -hydroxymethyl- β -hydroxyethyl)hydrazide, m. 130-5°, oxalic acid bis[2-[2-hydroxy-1-(hydroxymethyl)ethyl]hydrazide], m. 175.5-6.5°, oxalic acid bis[2-acetyl-2[2-acetoxy-1-(acetoxymethyl)ethyl]hydrazide], m. 187-8°; oxalic acid bis[2-(2-hydroxy-1-methylpropyl)hydrazide], m. 153.5-5.0°, oxalic acid hydrazide 2-[2-(benzylcarbamoylethyl)hydrazide], m. 244-5°, oxalic acid bis[2-[2-(benzylcarbamoylethyl)hydrazide], m. 216.5-18.0°, were also prepared. Acetic acid isopropylidenehydrazide (35.7 g.) in 300 ml. EtOH was hydrogenated at 50° in the presence of 1 g. PtO₂. The solution was evaporated to dryness and crystallized from C₆H₆-petr. ether to give 24.5 g. acetic acid 2-isopropylhydrazide, m. 53.5-5.0°. Also prepared were isobutyric acid 2-isopropylhydrazide, m. 69-70°, malonic acid bis(2-isopropylhydrazide), m. 127-9°, succinic acid bis(2-isopropylhydrazide), m. 158-9.5°, 2,2'-diisopropylcarbohydrazide, m. 141-2°, salicylic acid 2-isopropylhydrazide, m. 113-14°, 3,4,5-trimethoxybenzoic acid 2-isopropylhydrazide, m. 146.5-8.0°, 3,4,5-trimethoxybenzoic acid 2-(1-hydroxymethyl-2-hydroxyethyl)benzhydrazide, m. 111-12.5°, 2-(1-hydroxymethylethyl)2-hydroxyisonicotinic acid hydrazide, m. 124-7°, 3-hydroxy-2-naphthoic acid 2-isopropylhydrazide, m. 146-7°, indole-3-acetic acid 2-isopropylhydrazide, m. 130-1°.

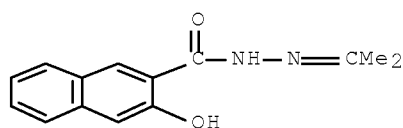
IT 73190-31-5P, 2-Naphthoic acid, 3-hydroxy-,
isopropylidenehydrazide 92255-70-4P, 2-Naphthoic acid,
3-hydroxy-, 2-isopropylhydrazide

RL: PREP (Preparation)

(preparation of)

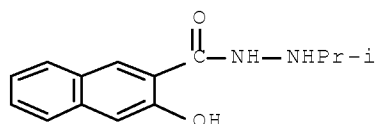
RN 73190-31-5 HCAPLUS

CN 2-Naphthalenecarboxylic acid, 3-hydroxy-, 2-(1-methylethylidene)hydrazide (CA INDEX NAME)



RN 92255-70-4 HCAPLUS

CN 2-Naphthalenecarboxylic acid, 3-hydroxy-, 2-(1-methylethyl)hydrazide
(CA INDEX NAME)



CC 26 (General Organic Chemistry)

IT Hydrazides

(amine oxidase inhibition by)

IT 3840-68-4P, Oxalic acid, bis(isopropylidenehydrazide) 5448-47-5P, Indole-3-acetic acid, hydrazide 6232-97-9P, Isobutyric acid, 2-isopropylhydrazide 7034-09-5P, Isonicotinic acid, [2-hydroxy-1-(hydroxymethyl)ethylidene]hydrazide 7034-12-0P, Isonicotinic acid, 2-[2-hydroxy-1-(hydroxymethyl)ethyl]hydrazide 13304-62-6P, Acrylamide, N-benzyl- 15316-21-9P, Indole-3-acetic acid, isopropylidenehydrazide 15317-49-4P, Indole-3-acetic acid, 2-isopropylhydrazide 15563-12-9P, Benzoic acid, 3,4,5-trimethoxy-, 2-isopropylhydrazide 18658-79-2P, Oxalic acid, bis(ethylidenehydrazide) 19411-38-2P, Oxalic acid, bis(sec-butylidenehydrazide) 21909-51-3P, Indole-3-acetic acid, 2-methyl-, hydrazide 26824-43-1P, Malonic acid, bis(2-isopropylhydrazide) 35532-19-5P, Oxalic acid, bis(2-ethylhydrazide) 35532-21-9P, Oxalic acid, bis[2-(2-hydroxy-1-methylpropyl)hydrazide] 35532-33-3P, Carbohydrazide, 1,5-diisopropyl- 63970-76-3P, Oxalic acid, methyl ester, hydrazide 73190-31-5P, 2-Naphthoic acid, 3-hydroxy-, isopropylidenehydrazide 83420-60-4P, Oxalic acid, bis(2-benzylhydrazide) 89620-31-5P, Oxalic acid, bis(cyclopentylidenehydrazide) 89620-45-1P, Oxalic acid, bis[(1-cyclopropylethylidene)hydrazide] 91346-66-6P, Cyclohexanecarboxylic acid, 2-hydroxy-, 2-isopropylhydrazide 91773-17-0P, Oxalic acid, 2-ethylhydrazide 2-isopropylhydrazide 91905-54-3P, Benzoic acid, 3,4,5-trimethoxy-, 2-[2-hydroxy-1-(hydroxymethyl)ethyl]hydrazide 92203-53-7P, Oxalic acid, bis(propylidenehydrazide) 92203-88-8P, Oxalic acid, ethylidenehydrazide 2-isopropylhydrazide 92223-99-9P, Oxalic acid, bis[[2-hydroxy-1-(hydroxymethyl)ethylidene]hydrazide] 92255-70-4P, 2-Naphthoic acid, 3-hydroxy-, 2-isopropylhydrazide 92351-54-7P, Oxalic acid, bis[2-[2-hydroxy-1-(hydroxymethyl)ethyl]hydrazide] 92351-56-9P, Oxalic acid, 2-[2-hydroxy-1-(hydroxymethyl)ethyl]hydrazide 2-isopropylhydrazide 92351-59-2P, Oxalic acid, bis(2-propylhydrazide) 92402-25-0P, Oxalic acid, bis(2-isopropylhydrazide) 93786-61-9P, Oxalic acid, bis(2-butylhydrazide) 93786-62-0P, Oxalic acid, bis(2-sec-butylhydrazide) 93786-63-1P, Oxalic acid,

bis(2-isobutylhydrazide) 94072-96-5P, Oxalic acid,
bis(butylidenehydrazide) 94072-97-6P, Oxalic acid,
bis(isobutylidenehydrazide) 94215-72-2P, Succinic acid,
bis(2-isopropylhydrazide) 94628-97-4P, Oxalic acid,
bis[(1-methylacetyliden)hydrazide] 94688-24-1P, Hydrazine,
1,2-bis(indol-3-ylglyoxyloyl)- 96130-96-0P, Oxalic acid,
2-[2-(benzylcarbamoyl)ethyl]hydrazide hydrazide 96417-86-6P,
Oxalic acid, bis(2-cyclopentylhydrazide) 96417-87-7P, Oxalic acid,
bis[2-(1-cyclopropylethyl)hydrazide] 96417-88-8P, Oxalic acid,
bis[(1-ethylpropyliden)hydrazide] 96591-42-3P, Benzoic acid,
3,4,5-trimethoxy-, [2-hydroxy-1-(hydroxymethyl)ethyliden]hydrazide
96874-51-0P, Oxalic acid, bis[2-(1-ethylpropyl)hydrazide]
97174-36-2P, Oxalic acid, 2-isopropylhydrazide 2-(α -
methylphenethyl)hydrazide 97281-74-8P, Oxalic acid,
2-isopropylhydrazide (α -methylphenethyliden)hydrazide
97491-38-8P, Oxalic acid, bis[[2-hydroxy-1-
(hydroxymethyl)ethyliden]hydrazide], tetraacetate 97528-04-6P,
Oxalic acid, bis[[2-hydroxy-1-(hydroxymethyl)ethyl]hydrazide],
tetraacetate diperchlorate 98636-67-0P, Oxalic acid,
bis(2-phenethylhydrazide) 99904-64-0P, Oxalic acid,
bis[(α -methylphenethyliden)hydrazide] 99998-68-2P, Oxalic
acid, bis[2-(α -methylphenethyl)hydrazide] 101122-35-4P,
Oxalic acid, bis[2-acetyl-2-[2-hydroxy-1-
(hydroxymethyl)ethyl]hydrazide], tetraacetate 101502-01-6P, Oxalic
acid, bis[2-[2-(benzylcarbamoyl)ethyl]hydrazide]
RL: PREP (Preparation)
(preparation of)

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